

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021721\
 Data File : BM028780.D
 Acq On : 17 Feb 2021 20:05
 Operator : CG/JU
 Sample : M1379-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBGY8

Manual Integrations
 APPROVED

mohammad
 2/18/2021 1:39:20 PM

Quant Time: Feb 18 04:46:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 18 04:11:41 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	490	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	1942	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	1098	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	2272	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	2684	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	2067	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	494	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1100	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	659	0.136	ng/ul#	83
5) 2-Methylnaphthalene	12.36	142	453	0.140	ng/ul	99
7) Acenaphthylene	14.26	152	1081	0.241	ng/ul#	92
8) Acenaphthene	14.60	153	1946	0.548	ng/ul	99
9) Fluorene	15.59	166	5880	1.491	ng/ul	97
11) Pentachlorophenol	16.94	266	19	0.032	ng/ul	94
12) Phenanthrene	17.31	178	51465	8.104	ng/ul	94
13) Anthracene	17.41	178	4493	0.742	ng/ul	97
15) Fluoranthene	19.32	202	68160	9.429	ng/ul	93
17) Pyrene	19.68	202	77839	7.982	ng/ul	95
18) Benzo(a)anthracene	21.41	228	8146	0.889	ng/ul	96
19) Chrysene	21.46	228	9970	1.095	ng/ul	95
21) Benzo(b)fluoranthene	22.98	252	8151m	1.152	ng/ul	
22) Benzo(k)fluoranthene	23.03	252	2746m	0.389	ng/ul	
23) Benzo(a)pyrene	23.56	252	2733	0.434	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.91	276	2363	0.288	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	673	0.098	ng/ul#	6
26) Benzo(g,h,i)perylene	26.60	276	2094	0.307	ng/ul	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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